



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:05 PM UTC

PDB ID : 6YV9 / pdb_00006yv9
Title : Mannosyltransferase PcManGT from *Pyrobaculum calidifontis* in complex with GDP-Man and Mn²⁺
Authors : Divne, C.; Gandini, R.
Deposited on : 2020-04-28
Resolution : 2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

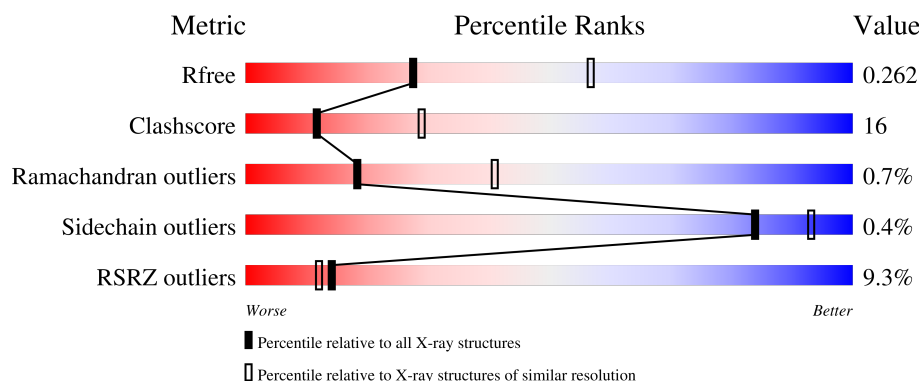
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

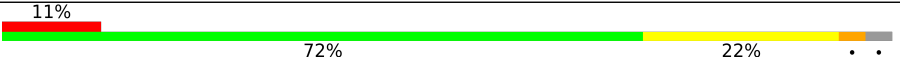
The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	355	
2	B	356	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5466 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

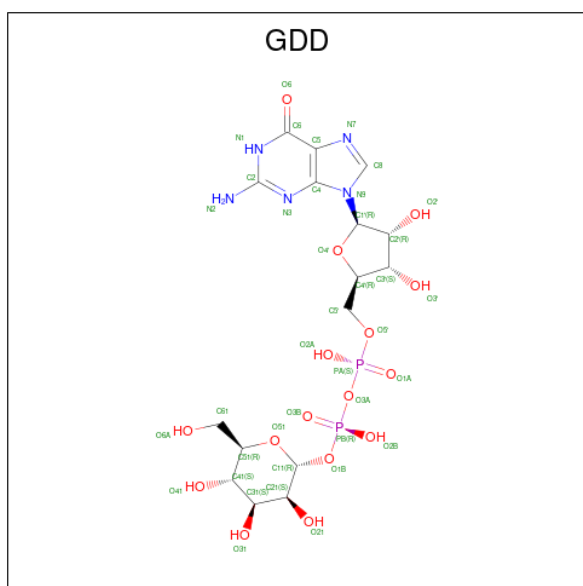
- Molecule 1 is a protein called Glycosyl transferase, family 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	345	Total	C	N	O	S	0	0	0
			2689	1734	476	474	5			

- Molecule 2 is a protein called Glycosyl transferase, family 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	346	Total	C	N	O	S	0	0	0
			2697	1739	477	475	6			

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE-ALPHA-D-MANNOSE (CCD ID: GDD) (formula: $C_{16}H_{25}N_5O_{16}P_2$).



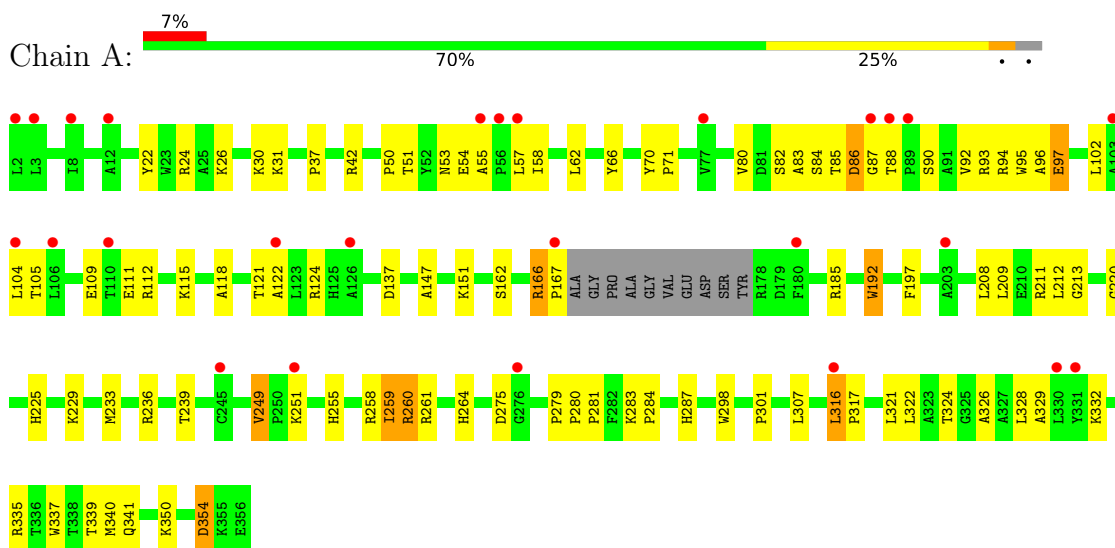
- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Mn 1	0	0
4	B	1	Total 1	Mn 1	0	0

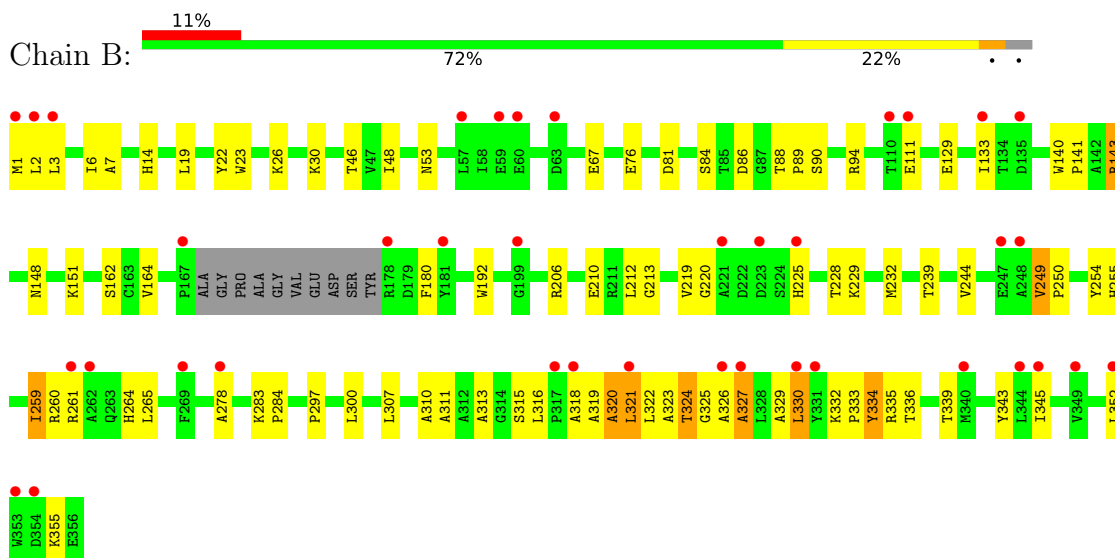
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Glycosyl transferase, family 2



• Molecule 2: Glycosyl transferase, family 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.47Å 107.42Å 163.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.41 – 2.70 46.41 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.41-2.70) 99.5 (46.41-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, R_{free}	0.215 , 0.265 0.217 , 0.262	Depositor DCC
R_{free} test set	2000 reflections (7.77%)	wwPDB-VP
Wilson B-factor (Å ²)	71.9	Xtriage
Anisotropy	0.397	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 65.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5466	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GDD, MN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.76	3/2766 (0.1%)	0.91	10/3789 (0.3%)
2	B	0.90	4/2774 (0.1%)	1.03	14/3799 (0.4%)
All	All	0.83	7/5540 (0.1%)	0.97	24/7588 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	B	0	3
All	All	0	4

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	334	TYR	C-N	30.34	1.71	1.33
1	A	259	ILE	C-N	21.42	1.63	1.33
2	B	327	ALA	C-N	16.71	1.57	1.33
1	A	260	ARG	C-N	15.37	1.56	1.33
2	B	261	ARG	C-N	9.60	1.46	1.33
1	A	166	ARG	CZ-NH1	5.59	1.40	1.32
2	B	260	ARG	C-N	-5.19	1.27	1.33

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	327	ALA	O-C-N	-19.48	101.77	122.03
2	B	249	VAL	CA-C-N	11.14	131.42	119.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	249	VAL	C-N-CA	11.14	131.42	119.05
1	A	249	VAL	CA-C-N	10.89	131.14	119.05
1	A	249	VAL	C-N-CA	10.89	131.14	119.05
2	B	143	ARG	CB-CG-CD	-10.18	87.89	111.30
2	B	129	GLU	CA-CB-CG	8.20	130.50	114.10
2	B	320	ALA	CA-C-N	-8.09	109.81	122.67
2	B	320	ALA	C-N-CA	-8.09	109.81	122.67
2	B	111	GLU	CA-CB-CG	7.54	129.19	114.10
1	A	259	ILE	O-C-N	-6.87	114.76	121.83
2	B	26	LYS	CG-CD-CE	-6.51	96.33	111.30
2	B	254	TYR	N-CA-C	6.41	117.93	111.07
2	B	324	THR	N-CA-CB	6.35	119.47	109.82
1	A	97	GLU	CA-CB-CG	6.23	126.57	114.10
1	A	259	ILE	CA-C-N	6.02	133.03	121.54
1	A	259	ILE	C-N-CA	6.02	133.03	121.54
1	A	166	ARG	NE-CZ-NH1	5.78	127.28	121.50
2	B	324	THR	CB-CA-C	-5.72	101.82	110.92
1	A	275	ASP	N-CA-C	5.35	117.14	108.26
2	B	129	GLU	N-CA-CB	5.15	118.94	110.39
2	B	261	ARG	O-C-N	5.07	127.50	122.03
1	A	97	GLU	CA-C-N	-5.03	111.78	121.58
1	A	97	GLU	C-N-CA	-5.03	111.78	121.58

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	97	GLU	Sidechain
2	B	321	LEU	Peptide
2	B	327	ALA	Mainchain
2	B	330	LEU	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2689	0	2704	98	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2697	0	2715	76	0
3	A	39	0	23	6	0
3	B	39	0	22	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
All	All	5466	0	5464	174	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (174) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:334:TYR:C	2:B:335:ARG:N	1.71	1.48
1:A:332:LYS:HD2	1:A:335:ARG:NH2	1.47	1.28
2:B:330:LEU:CD2	2:B:335:ARG:HD2	1.77	1.14
1:A:332:LYS:CD	1:A:335:ARG:NH2	2.22	1.03
2:B:330:LEU:HD21	2:B:335:ARG:HD2	1.39	1.02
1:A:332:LYS:HD2	1:A:335:ARG:CZ	1.96	0.95
1:A:332:LYS:HD2	1:A:335:ARG:HH21	1.14	0.93
2:B:1:MET:H2	2:B:3:LEU:HD23	1.33	0.91
1:A:192:TRP:HA	1:A:236:ARG:NH1	1.90	0.87
1:A:83:ALA:HB3	1:A:111:GLU:HA	1.56	0.86
2:B:321:LEU:O	2:B:324:THR:OG1	1.94	0.84
2:B:330:LEU:CD2	2:B:335:ARG:CD	2.56	0.84
1:A:82:SER:HA	1:A:109:GLU:HB2	1.61	0.82
2:B:212:LEU:HD22	2:B:229:LYS:HE3	1.60	0.82
1:A:54:GLU:H	1:A:88:THR:HG21	1.45	0.82
1:A:93:ARG:HB2	1:A:93:ARG:NH1	1.94	0.81
1:A:332:LYS:CE	1:A:335:ARG:NH2	2.43	0.80
2:B:1:MET:N	2:B:3:LEU:HD23	1.95	0.80
1:A:332:LYS:CD	1:A:335:ARG:CZ	2.58	0.79
2:B:326:ALA:O	2:B:329:ALA:HB3	1.82	0.78
1:A:332:LYS:HE3	1:A:335:ARG:CZ	2.15	0.77
2:B:330:LEU:HD21	2:B:335:ARG:CD	2.17	0.74
2:B:90:SER:O	2:B:94:ARG:HG3	1.87	0.74
2:B:330:LEU:HD23	2:B:335:ARG:HD2	1.68	0.73
1:A:332:LYS:HE3	1:A:335:ARG:NH2	2.02	0.73
1:A:307:LEU:HB3	1:A:322:LEU:HD11	1.71	0.73
1:A:166:ARG:HB2	1:A:167:PRO:HD2	1.72	0.72
1:A:220:GLY:HA3	1:A:264:HIS:CD2	2.25	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:311:ALA:HA	2:B:316:LEU:HD21	1.73	0.70
1:A:80:VAL:HG21	1:A:122:ALA:HB2	1.74	0.69
1:A:53:ASN:OD1	1:A:84:SER:OG	2.10	0.69
2:B:67:GLU:O	2:B:143:ARG:HB3	1.93	0.68
1:A:93:ARG:HB2	1:A:93:ARG:HH11	1.56	0.68
1:A:316:LEU:H	1:A:317:PRO:HD2	1.59	0.68
2:B:321:LEU:HA	2:B:324:THR:H	1.57	0.67
2:B:325:GLY:O	2:B:329:ALA:HB2	1.94	0.67
1:A:51:THR:HG21	1:A:58:ILE:HD13	1.76	0.66
1:A:42:ARG:HD2	1:A:71:PRO:HG3	1.78	0.65
1:A:192:TRP:HA	1:A:236:ARG:HH11	1.61	0.65
1:A:258:ARG:NH2	1:A:340:MET:SD	2.69	0.64
2:B:255:HIS:O	2:B:259:ILE:HG12	1.96	0.64
1:A:55:ALA:O	1:A:58:ILE:HG22	1.97	0.64
2:B:30:LYS:O	2:B:30:LYS:HD3	1.97	0.64
1:A:261:ARG:CZ	3:A:400:GDD:H51	2.29	0.63
2:B:311:ALA:HA	2:B:316:LEU:CD2	2.28	0.63
2:B:321:LEU:HG	2:B:324:THR:HG23	1.81	0.63
1:A:332:LYS:CE	1:A:335:ARG:CZ	2.76	0.62
2:B:320:ALA:O	2:B:321:LEU:HD12	2.00	0.62
2:B:278:ALA:HB3	2:B:283:LYS:HD3	1.82	0.60
1:A:66:TYR:HB2	1:A:95:TRP:HZ2	1.67	0.60
1:A:24:ARG:HB3	1:A:24:ARG:NH1	2.16	0.59
1:A:93:ARG:HB2	1:A:93:ARG:CZ	2.32	0.59
1:A:192:TRP:HA	1:A:236:ARG:HH12	1.68	0.59
2:B:311:ALA:HB1	2:B:319:ALA:HB2	1.83	0.59
2:B:311:ALA:O	2:B:316:LEU:HD21	2.03	0.58
2:B:307:LEU:HD22	2:B:318:ALA:HB1	1.86	0.57
1:A:279:PRO:O	1:A:280:PRO:C	2.45	0.57
1:A:58:ILE:O	1:A:62:LEU:HD12	2.05	0.57
2:B:320:ALA:O	2:B:323:ALA:HB3	2.05	0.57
1:A:332:LYS:HD2	1:A:335:ARG:NE	2.20	0.56
1:A:115:LYS:NZ	3:A:400:GDD:O21	2.28	0.56
2:B:219:VAL:HG11	2:B:225:HIS:CG	2.40	0.56
1:A:104:LEU:HD23	1:A:105:THR:N	2.21	0.56
2:B:316:LEU:HD22	2:B:319:ALA:CB	2.34	0.56
1:A:95:TRP:HH2	1:A:102:LEU:HD23	1.71	0.56
1:A:335:ARG:O	1:A:339:THR:HG23	2.06	0.55
1:A:279:PRO:C	1:A:281:PRO:HD2	2.32	0.55
2:B:162:SER:HB3	2:B:239:THR:HG22	1.89	0.55
1:A:249:VAL:HG13	1:A:249:VAL:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1:MET:HG3	2:B:2:LEU:H	1.72	0.54
2:B:311:ALA:CA	2:B:316:LEU:HD21	2.37	0.54
1:A:283:LYS:HE3	1:A:287:HIS:CD2	2.43	0.54
1:A:86:ASP:OD1	1:A:88:THR:HG23	2.07	0.54
2:B:320:ALA:C	2:B:321:LEU:HD12	2.33	0.54
1:A:37:PRO:HB2	1:A:151:LYS:HE3	1.90	0.54
1:A:147:ALA:O	1:A:151:LYS:HG3	2.08	0.54
1:A:30:LYS:O	1:A:30:LYS:HD2	2.07	0.53
1:A:90:SER:O	1:A:94:ARG:NE	2.42	0.53
3:A:400:GDD:O5'	3:A:400:GDD:H8	2.09	0.53
1:A:83:ALA:HB2	1:A:109:GLU:O	2.09	0.53
1:A:212:LEU:N	1:A:213:GLY:HA2	2.23	0.53
1:A:298:TRP:O	1:A:301:PRO:HD2	2.10	0.52
1:A:316:LEU:N	1:A:317:PRO:HD2	2.24	0.52
1:A:192:TRP:CA	1:A:236:ARG:NH1	2.68	0.51
2:B:329:ALA:HA	2:B:334:TYR:CD2	2.46	0.51
2:B:319:ALA:HA	2:B:322:LEU:HB3	1.91	0.51
1:A:260:ARG:NH1	3:A:400:GDD:O3B	2.43	0.51
2:B:315:SER:O	2:B:316:LEU:HD23	2.11	0.50
2:B:148:ASN:O	2:B:151:LYS:HG2	2.11	0.50
1:A:85:THR:O	1:A:87:GLY:N	2.42	0.50
1:A:57:LEU:HD13	1:A:137:ASP:HB3	1.94	0.50
1:A:162:SER:HB3	1:A:239:THR:HG22	1.93	0.49
1:A:121:THR:O	1:A:124:ARG:HG2	2.11	0.49
1:A:307:LEU:HB3	1:A:322:LEU:CD1	2.41	0.49
1:A:324:THR:O	1:A:328:LEU:HD13	2.12	0.49
2:B:352:LEU:O	2:B:355:LYS:HG2	2.11	0.49
2:B:1:MET:H2	2:B:3:LEU:CD2	2.17	0.49
1:A:90:SER:O	1:A:94:ARG:HG2	2.12	0.48
2:B:7:ALA:HB1	2:B:307:LEU:HG	1.95	0.48
2:B:206:ARG:NH1	2:B:210:GLU:OE2	2.46	0.48
2:B:212:LEU:N	2:B:213:GLY:HA2	2.28	0.48
2:B:3:LEU:HG	2:B:310:ALA:HB2	1.95	0.48
2:B:228:THR:O	2:B:232:MET:HG2	2.12	0.48
1:A:82:SER:HA	1:A:109:GLU:CB	2.40	0.48
2:B:81:ASP:OD1	2:B:84:SER:OG	2.28	0.47
1:A:95:TRP:CH2	1:A:102:LEU:HD23	2.50	0.47
2:B:283:LYS:N	2:B:284:PRO:HD2	2.29	0.47
1:A:208:LEU:HD22	1:A:233:MET:HE1	1.97	0.47
1:A:212:LEU:H	1:A:213:GLY:HA2	1.79	0.47
1:A:261:ARG:HG3	1:A:261:ARG:NH1	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:249:VAL:HA	2:B:250:PRO:HD3	1.55	0.47
2:B:2:LEU:O	2:B:6:ILE:HG12	2.15	0.46
1:A:192:TRP:O	1:A:236:ARG:HG2	2.15	0.46
2:B:19:LEU:HD21	2:B:23:TRP:CZ2	2.51	0.46
2:B:316:LEU:HD22	2:B:319:ALA:HB3	1.98	0.46
1:A:329:ALA:O	1:A:335:ARG:HD3	2.16	0.46
2:B:332:LYS:N	2:B:333:PRO:HD2	2.30	0.46
2:B:316:LEU:HD23	2:B:316:LEU:HA	1.45	0.46
1:A:54:GLU:N	1:A:88:THR:HG21	2.22	0.46
1:A:332:LYS:CD	1:A:335:ARG:NE	2.79	0.45
2:B:53:ASN:HA	2:B:86:ASP:OD2	2.16	0.45
1:A:31:LYS:NZ	1:A:31:LYS:HB2	2.31	0.45
2:B:297:PRO:O	2:B:345:ILE:HD11	2.16	0.45
2:B:14:HIS:CD2	2:B:300:LEU:HD13	2.52	0.45
1:A:350:LYS:O	1:A:354:ASP:HB2	2.16	0.45
2:B:212:LEU:HD23	2:B:212:LEU:HA	1.74	0.45
1:A:332:LYS:HE3	1:A:335:ARG:NH1	2.31	0.45
2:B:46:THR:HG23	2:B:76:GLU:HG3	1.99	0.45
1:A:287:HIS:CD2	2:B:313:ALA:HB2	2.52	0.45
2:B:220:GLY:HA3	2:B:264:HIS:ND1	2.32	0.44
2:B:88:THR:HB	2:B:89:PRO:HD3	2.00	0.44
2:B:48:ILE:HB	2:B:133:ILE:HD13	2.00	0.44
2:B:326:ALA:O	2:B:329:ALA:CB	2.59	0.44
1:A:93:ARG:HA	1:A:96:ALA:HB3	1.99	0.44
1:A:22:TYR:CE2	1:A:26:LYS:HD3	2.52	0.44
1:A:24:ARG:HB3	1:A:24:ARG:HH11	1.80	0.44
2:B:164:VAL:O	2:B:244:VAL:HA	2.18	0.44
1:A:93:ARG:HH11	1:A:93:ARG:CB	2.28	0.44
1:A:109:GLU:OE2	1:A:118:ALA:HA	2.18	0.43
1:A:211:ARG:O	1:A:211:ARG:HG2	2.18	0.43
2:B:22:TYR:HB2	2:B:180:PHE:HE1	1.83	0.43
1:A:70:TYR:CG	1:A:71:PRO:HD2	2.54	0.43
1:A:112:ARG:HG3	3:A:400:GDD:O6	2.19	0.43
1:A:317:PRO:O	1:A:321:LEU:HG	2.18	0.43
2:B:140:TRP:HA	2:B:141:PRO:HD3	1.89	0.43
2:B:311:ALA:C	2:B:316:LEU:HD21	2.44	0.43
2:B:316:LEU:HD22	2:B:319:ALA:HB2	2.01	0.43
2:B:307:LEU:HD23	2:B:307:LEU:HA	1.69	0.42
1:A:51:THR:HA	3:A:400:GDD:O2'	2.19	0.42
1:A:255:HIS:O	1:A:259:ILE:HG13	2.19	0.42
1:A:55:ALA:HB2	1:A:88:THR:HG22	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:86:ASP:OD2	2:B:86:ASP:N	2.50	0.42
1:A:316:LEU:H	1:A:317:PRO:CD	2.31	0.41
2:B:339:THR:HG22	2:B:343:TYR:CE2	2.54	0.41
1:A:337:TRP:O	1:A:341:GLN:HG2	2.20	0.41
1:A:209:LEU:O	1:A:213:GLY:HA2	2.21	0.41
1:A:85:THR:OG1	1:A:86:ASP:N	2.52	0.41
1:A:332:LYS:CD	1:A:335:ARG:HH21	2.02	0.41
2:B:14:HIS:CE1	2:B:300:LEU:HB2	2.56	0.41
1:A:162:SER:O	1:A:239:THR:HA	2.21	0.41
1:A:192:TRP:CA	1:A:236:ARG:HH12	2.31	0.41
1:A:225:HIS:NE2	1:A:229:LYS:HE2	2.35	0.41
2:B:249:VAL:O	2:B:249:VAL:HG23	2.20	0.41
2:B:332:LYS:O	2:B:336:THR:HG23	2.20	0.41
1:A:50:PRO:HG3	1:A:115:LYS:HE3	2.03	0.40
1:A:185:ARG:NE	1:A:185:ARG:HA	2.35	0.40
1:A:162:SER:HB2	1:A:197:PHE:CD2	2.57	0.40
1:A:283:LYS:N	1:A:284:PRO:HD2	2.37	0.40
2:B:30:LYS:HD3	2:B:30:LYS:C	2.46	0.40
2:B:265:LEU:HD23	2:B:265:LEU:HA	1.85	0.40
2:B:329:ALA:HA	2:B:334:TYR:CE2	2.56	0.40
1:A:326:ALA:O	1:A:329:ALA:HB3	2.21	0.40
2:B:259:ILE:HG12	2:B:259:ILE:H	1.68	0.40
1:A:88:THR:O	1:A:92:VAL:HG23	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	341/355 (96%)	327 (96%)	10 (3%)	4 (1%)	10	27
2	B	342/356 (96%)	328 (96%)	13 (4%)	1 (0%)	36	60

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	683/711 (96%)	655 (96%)	23 (3%)	5 (1%)	18	41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	251	LYS
1	A	192	TRP
1	A	86	ASP
2	B	192	TRP
1	A	316	LEU

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	267/273 (98%)	266 (100%)	1 (0%)	84	93
2	B	268/274 (98%)	267 (100%)	1 (0%)	84	93
All	All	535/547 (98%)	533 (100%)	2 (0%)	84	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	354	ASP
2	B	259	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	117	HIS
1	A	125	HIS
1	A	264	HIS
1	A	287	HIS
2	B	125	HIS
2	B	182	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GDD	B	400	-	41,42,42	1.90	13 (31%)	62,65,65	1.68	15 (24%)
3	GDD	A	400	4	41,42,42	1.88	12 (29%)	62,65,65	1.65	16 (25%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GDD	B	400	-	-	5/23/59/59	0/4/4/4
3	GDD	A	400	4	-	3/23/59/59	0/4/4/4

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	400	GDD	C4-N3	5.95	1.47	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	400	GDD	C4-N3	5.89	1.47	1.34
3	A	400	GDD	C2-N2	4.02	1.43	1.34
3	B	400	GDD	C2-N2	3.88	1.43	1.34
3	A	400	GDD	C3'-C4'	-3.41	1.44	1.53
3	B	400	GDD	C3'-C4'	-3.37	1.44	1.53
3	B	400	GDD	C2-N1	3.17	1.45	1.37
3	A	400	GDD	C2-N1	3.12	1.45	1.37
3	A	400	GDD	C3'-C2'	-3.05	1.45	1.53
3	B	400	GDD	O4'-C1'	2.92	1.48	1.42
3	B	400	GDD	O21-C21	-2.87	1.35	1.43
3	A	400	GDD	C2'-C1'	-2.72	1.44	1.53
3	B	400	GDD	C2'-C1'	-2.61	1.45	1.53
3	A	400	GDD	O21-C21	-2.46	1.36	1.43
3	B	400	GDD	C3'-C2'	-2.41	1.46	1.53
3	A	400	GDD	O4'-C1'	2.37	1.47	1.42
3	A	400	GDD	C5'-C4'	-2.37	1.44	1.51
3	A	400	GDD	C5-N7	-2.32	1.34	1.39
3	A	400	GDD	C1'-N9	-2.32	1.41	1.47
3	B	400	GDD	C5'-C4'	-2.30	1.44	1.51
3	B	400	GDD	C5-N7	-2.24	1.34	1.39
3	B	400	GDD	C41-C51	-2.23	1.48	1.53
3	B	400	GDD	O31-C31	-2.16	1.37	1.43
3	B	400	GDD	O4'-C4'	2.12	1.49	1.45
3	A	400	GDD	O4'-C4'	2.08	1.49	1.45

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	400	GDD	C5-C4-N3	-4.52	121.19	128.39
3	A	400	GDD	C2-N3-C4	4.34	119.78	112.30
3	A	400	GDD	C5-C4-N3	-4.34	121.48	128.39
3	B	400	GDD	C2-N3-C4	4.12	119.39	112.30
3	B	400	GDD	O51-C11-O1B	3.75	116.27	111.36
3	A	400	GDD	O1B-C11-C21	3.35	114.52	108.38
3	A	400	GDD	N9-C4-N3	3.28	132.51	125.95
3	B	400	GDD	C2'-C3'-C4'	3.11	108.61	102.61
3	B	400	GDD	C8-N7-C5	3.07	109.73	104.26
3	A	400	GDD	C31-C41-C51	3.06	115.79	110.23
3	A	400	GDD	O51-C51-C41	2.98	115.06	109.70
3	A	400	GDD	O6-C6-C5	-2.91	118.86	126.53
3	B	400	GDD	N9-C4-N3	2.85	131.66	125.95
3	B	400	GDD	C2-N1-C6	-2.83	119.97	125.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	400	GDD	C4'-O4'-C1'	-2.72	103.47	109.47
3	B	400	GDD	C5-C6-N1	2.70	120.13	113.25
3	B	400	GDD	C6-C5-N7	2.67	135.15	130.29
3	A	400	GDD	C8-N7-C5	2.56	108.83	104.26
3	B	400	GDD	N9-C8-N7	-2.52	108.72	113.40
3	B	400	GDD	C4-C5-N7	-2.50	106.72	110.67
3	A	400	GDD	C5-C6-N1	2.49	119.60	113.25
3	A	400	GDD	N9-C8-N7	-2.48	108.81	113.40
3	B	400	GDD	O6-C6-C5	-2.43	120.11	126.53
3	A	400	GDD	O2A-PA-O3A	2.36	113.66	107.27
3	B	400	GDD	O51-C51-C41	2.13	113.53	109.70
3	A	400	GDD	C2-N1-C6	-2.12	121.27	125.11
3	A	400	GDD	N2-C2-N1	2.11	121.22	116.76
3	A	400	GDD	C6-C5-N7	2.09	134.09	130.29
3	B	400	GDD	N2-C2-N1	2.05	121.09	116.76
3	B	400	GDD	O2A-PA-O3A	2.05	112.80	107.27
3	A	400	GDD	C61-C51-C41	-2.03	108.03	113.02

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	400	GDD	O51-C11-O1B-PB
3	B	400	GDD	C41-C51-C61-O6A
3	B	400	GDD	O51-C51-C61-O6A
3	A	400	GDD	PB-O3A-PA-O2A
3	B	400	GDD	PB-O3A-PA-O2A
3	B	400	GDD	O51-C11-O1B-PB
3	B	400	GDD	PB-O3A-PA-O1A
3	A	400	GDD	PB-O3A-PA-O1A

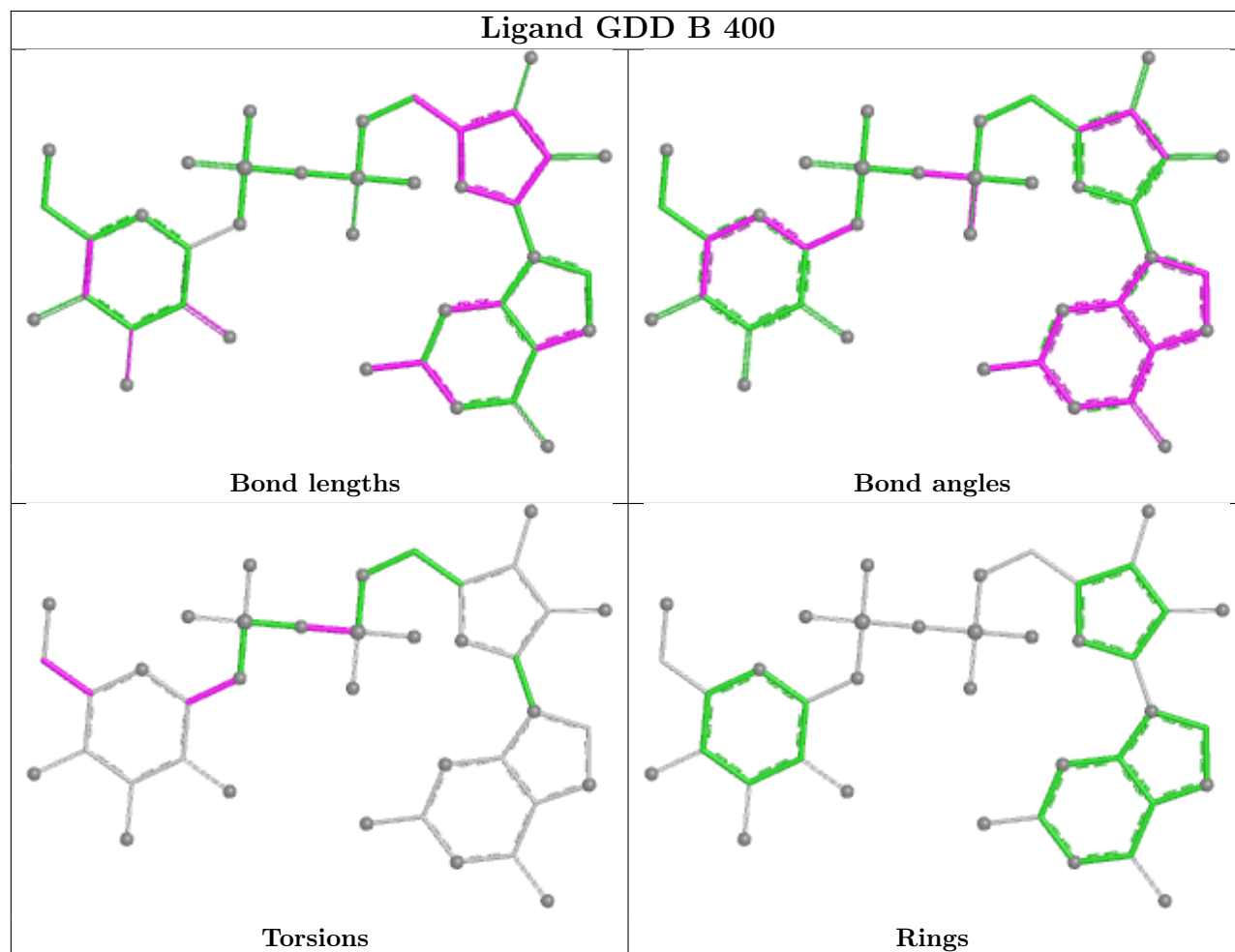
There are no ring outliers.

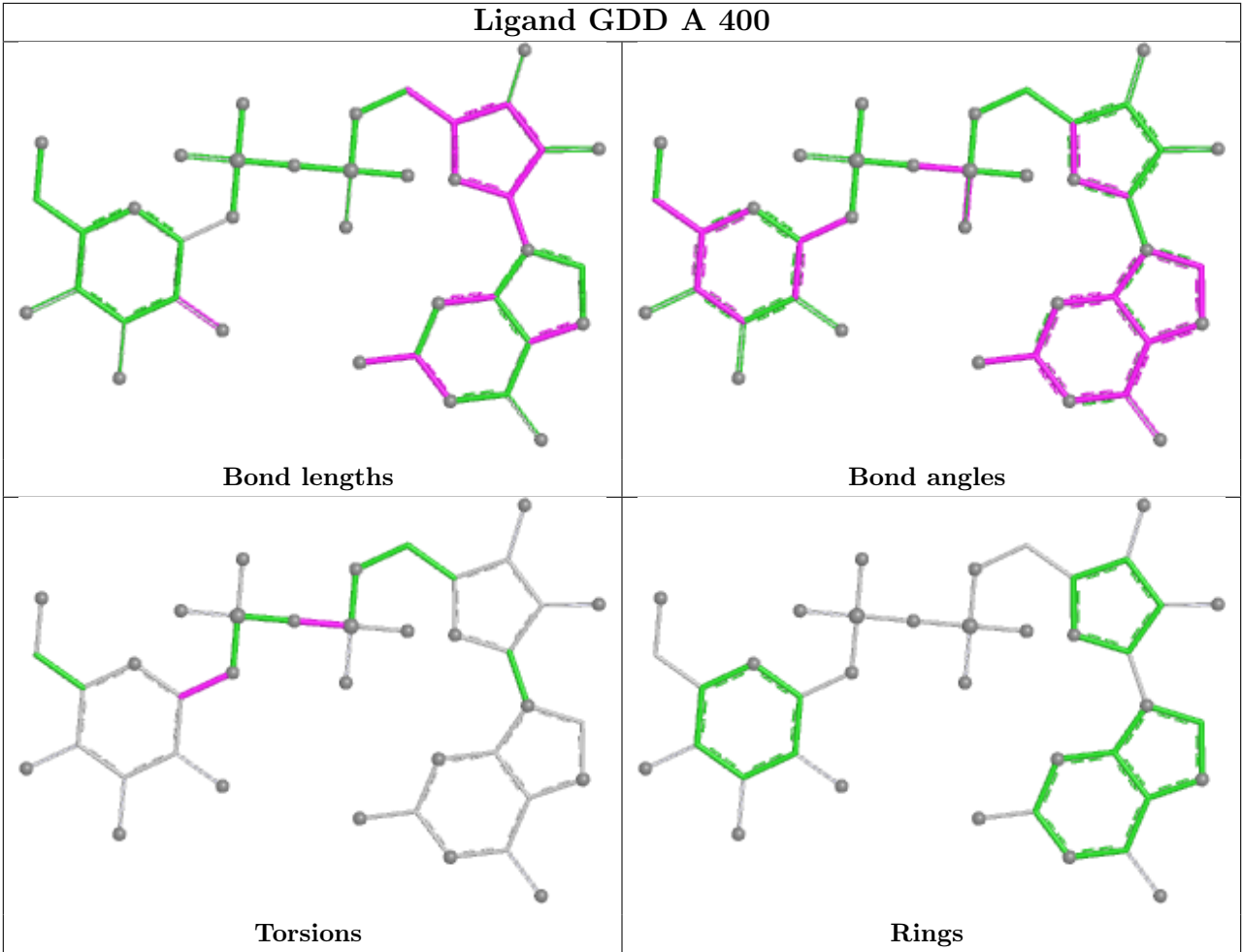
1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	400	GDD	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.

Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	B	1
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	334:TYR	C	335:ARG	N	1.71
1	A	259:ILE	C	260:ARG	N	1.63

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	345/355 (97%)	0.48	26 (7%) 20 17	39, 75, 122, 150	0
2	B	346/356 (97%)	0.47	38 (10%) 10 9	37, 69, 123, 158	0
All	All	691/711 (97%)	0.47	64 (9%) 14 12	37, 72, 123, 158	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	223	ASP	5.1
1	A	331	TYR	4.4
2	B	353	TRP	3.6
1	A	55	ALA	3.6
1	A	126	ALA	3.5
2	B	60	GLU	3.4
1	A	316	LEU	3.4
1	A	12	ALA	3.4
2	B	330	LEU	3.4
2	B	221	ALA	3.3
2	B	344	LEU	3.2
2	B	349	VAL	3.2
1	A	103	ALA	3.1
2	B	2	LEU	3.0
2	B	247	GLU	3.0
1	A	77	VAL	3.0
2	B	278	ALA	3.0
1	A	167	PRO	2.9
1	A	88	THR	2.9
1	A	57	LEU	2.9
1	A	104	LEU	2.9
1	A	245	CYS	2.9
2	B	135	ASP	2.9
2	B	167	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	251	LYS	2.8
2	B	248	ALA	2.8
2	B	262	ALA	2.8
2	B	331	TYR	2.8
2	B	345	ILE	2.7
2	B	3	LEU	2.7
2	B	317	PRO	2.6
2	B	327	ALA	2.6
1	A	122	ALA	2.5
1	A	180	PHE	2.5
2	B	57	LEU	2.5
1	A	56	PRO	2.5
1	A	330	LEU	2.5
2	B	225	HIS	2.5
2	B	63	ASP	2.4
2	B	133	ILE	2.4
2	B	59	GLU	2.4
1	A	110	THR	2.4
2	B	110	THR	2.3
1	A	276	GLY	2.3
2	B	321	LEU	2.3
2	B	111	GLU	2.3
2	B	340	MET	2.3
1	A	203	ALA	2.3
2	B	199	GLY	2.3
2	B	181	TYR	2.2
2	B	261	ARG	2.2
2	B	1	MET	2.2
2	B	318	ALA	2.2
2	B	352	LEU	2.2
2	B	269	PHE	2.2
1	A	89	PRO	2.2
1	A	2	LEU	2.2
1	A	3	LEU	2.2
1	A	87	GLY	2.1
2	B	178	ARG	2.1
1	A	106	LEU	2.1
2	B	326	ALA	2.1
1	A	8	ILE	2.1
2	B	354	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

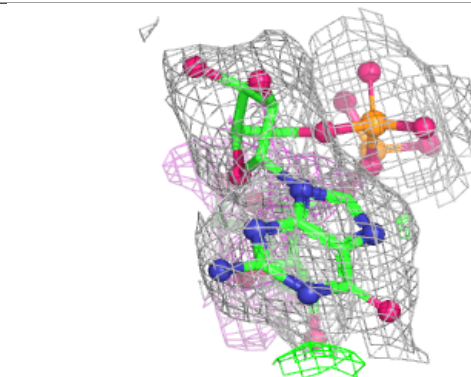
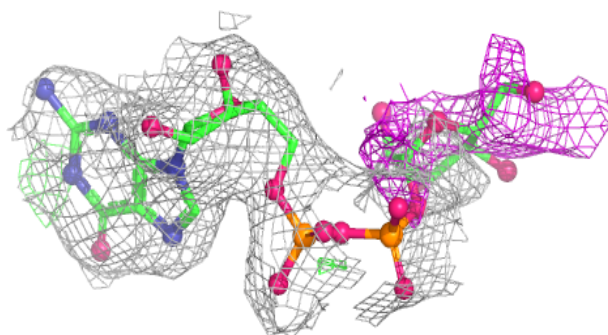
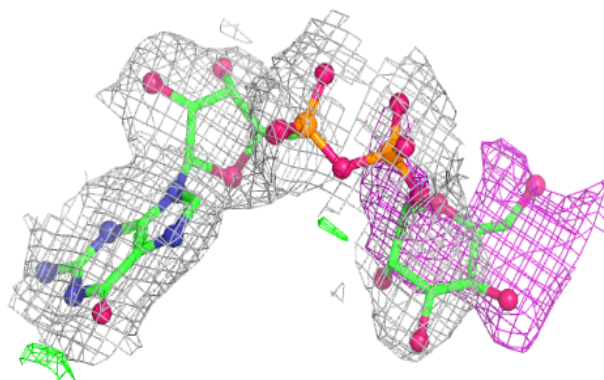
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GDD	B	400	39/39	0.86	0.12	51,82,126,134	0
3	GDD	A	400	39/39	0.87	0.10	67,94,126,127	0
4	MN	A	401	1/1	0.92	0.07	125,125,125,125	0
4	MN	B	401	1/1	0.93	0.06	129,129,129,129	0

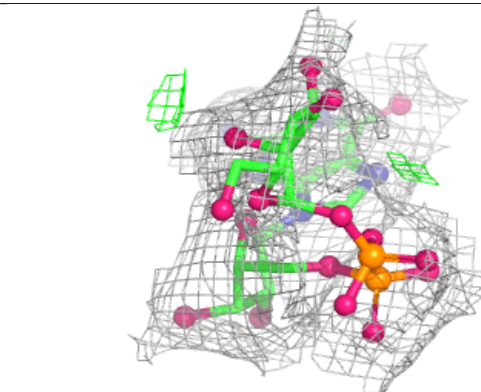
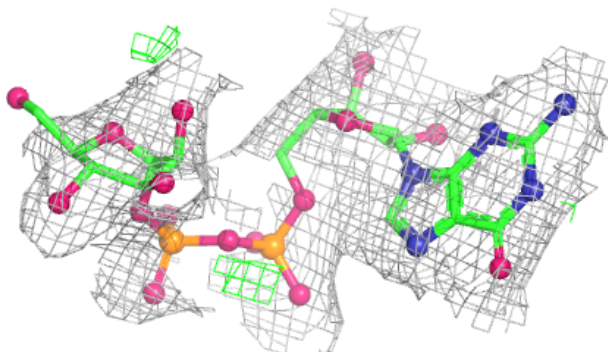
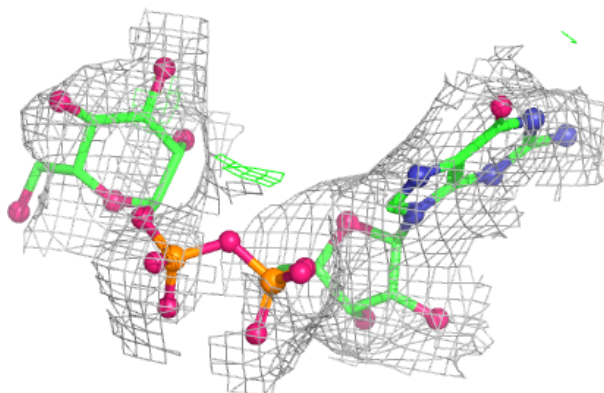
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around GDD B 400:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around GDD A 400:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.